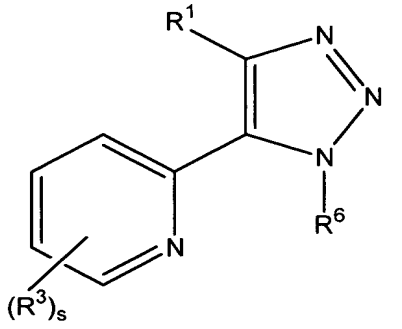
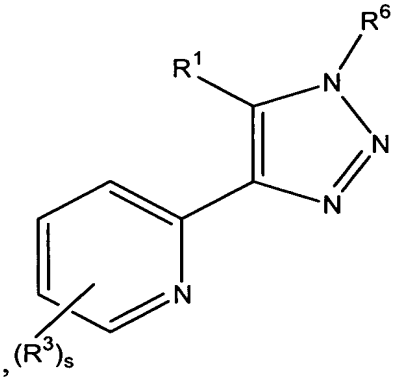
Amendment to the Claims

The claimed invention is:

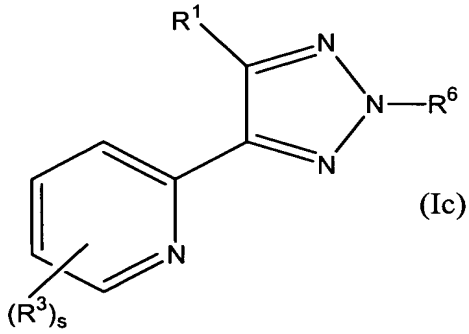
1. (Currently Amended) A compound of formula (Ia), (Ib), or (Ic):



(Ia)



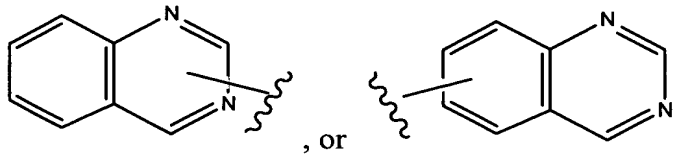
(Ib)



(Ic)

or a pharmaceutically acceptable salt, ~~prodrug~~, tautomer, hydrate or ~~solvate~~ thereof, wherein:

R^1 is a group of the formula



, or

wherein R^1 can optionally be further independently substituted with at least one moiety independently selected from the group consisting of: carbonyl, halo, halo(C_1-C_6)alkyl, perhalo(C_1-C_6)alkyl, perhalo(C_1-C_6)alkoxy, (C_1-C_6)alkyl, (C_2-C_6)alkenyl, (C_2-C_6)alkynyl, hydroxy, oxo, mercapto, (C_1-C_6)alkylthio, (C_1-C_6)alkoxy, (C_5-C_{10})aryl ~~or (C_5-C_{10})heteroaryl~~, (C_5-C_{10})aryloxy ~~or (C_5-C_{10})heteroaryloxy~~, (C_5-C_{10})ar(C_1-C_6)alkyl ~~or (C_5-C_{10})heteroar(C_1-C_6)alkyl~~,

$(C_5\text{-}C_{10})$ar$(C_1\text{-}C_6)$alkoxy ~~or $(C_5\text{-}C_{10})$heteroar$(C_1\text{-}C_6)$alkoxy~~, HO-(C=O)-, ester, amido, ether, amino, amino$(C_1\text{-}C_6)$alkyl, $(C_1\text{-}C_6)$alkylamino$(C_1\text{-}C_6)$alkyl, di$(C_1\text{-}C_6)$alkylamino$(C_1\text{-}C_6)$alkyl, ~~$(C_5\text{-}C_{10})$heterocyclyl$(C_1\text{-}C_6)$alkyl~~, $(C_1\text{-}C_6)$alkyl- and di$(C_1\text{-}C_6)$alkylamino, cyano, nitro, carbamoyl, $(C_1\text{-}C_6)$alkylcarbonyl, $(C_1\text{-}C_6)$alkoxycarbonyl, $(C_1\text{-}C_6)$alkylaminocarbonyl, di$(C_1\text{-}C_6)$alkylaminocarbonyl, $(C_5\text{-}C_{10})$arylcarbonyl, $(C_5\text{-}C_{10})$aryloxycarbonyl, $(C_1\text{-}C_6)$alkylsulfonyl, and $(C_5\text{-}C_{10})$arylsulfonyl;

each R^3 is independently selected from the group consisting of: hydrogen, halo, halo$(C_1\text{-}C_6)$alkyl, $(C_1\text{-}C_6)$alkyl, $(C_2\text{-}C_6)$alkenyl, $(C_2\text{-}C_6)$alkynyl, perhalo$(C_1\text{-}C_6)$alkyl, phenyl, $(C_3\text{-}C_{10})$cycloalkyl, hydroxy, $(C_1\text{-}C_6)$alkoxy, perhalo$(C_1\text{-}C_6)$alkoxy, phenoxy, $(C_3\text{-}C_{10})$cycloalkyl-O-, $(C_1\text{-}C_6)$alkyl-S-, $(C_1\text{-}C_6)$alkyl-SO_2-, $(C_1\text{-}C_6)$alkyl-NH-SO_2-, O_2N-, NC-, amino, Ph$(CH_2)_{1\text{-}6}$HN-, $(C_1\text{-}C_6)$alkyl HN-, $(C_1\text{-}C_6)$alkylamino, $[(C_1\text{-}C_6)\text{alkyl}]_2$-amino, $(C_1\text{-}C_6)$alkyl-SO_2-NH-, amino(C=O)-, aminoO_2S-, $(C_1\text{-}C_6)$alkyl-(C=O)-NH-, $(C_1\text{-}C_6)$alkyl-(C=O)-[(($(C_1\text{-}C_6)$alkyl)-N]-, phenyl-(C=O)-NH-, phenyl-(C=O)-[($(C_1\text{-}C_6)$alkyl)-N]-, $(C_1\text{-}C_6)$alkyl-(C=O)-, phenyl-(C=O)-, $(C_3\text{-}C_{10})$cycloalkyl-(C=O)-, HO-(C=O)-, $(C_1\text{-}C_6)$alkyl-O-(C=O)-, H_2N(C=O)-, $(C_1\text{-}C_6)$alkyl-NH-(C=O)-, $[(C_1\text{-}C_6)\text{alkyl}]_2$-N-(C=O)-, phenyl-NH-(C=O)-, phenyl-[(($C_1\text{-}C_6$)alkyl)-N]-(C=O)-, $(C_3\text{-}C_{10})$cycloalkyl-NH-(C=O)- and $(C_1\text{-}C_6)$alkyl-(C=O)-O-;

where alkyl, alkenyl, alkynyl, phenyl, cycloalkyl, alkoxy, phenoxy, amino of R^3 is optionally substituted by at least one substituent independently selected from $(C_1\text{-}C_6)$alkyl, $(C_1\text{-}C_6)$alkoxy, halo$(C_1\text{-}C_6)$alkyl, halo, H_2N-, Ph$(CH_2)_{1\text{-}6}$HN-, and $(C_1\text{-}C_6)$alkylHN-;

s is an integer from one to five;

and

R^6 is selected from the group consisting of hydrogen, $(C_1\text{-}C_6)$alkyl, $(C_2\text{-}C_6)$alkenyl, $(C_2\text{-}C_6)$alkynyl, phenyl, $(C_3\text{-}C_{10})$cycloalkyl, $(C_1\text{-}C_6)$alkyl-(SO_2)-, phenyl-(SO_2)-, H_2N-(SO_2)-, $(C_1\text{-}C_6)$alkyl-NH-(SO_2)-, $((C_1\text{-}C_6)\text{alkyl})_2$N-$(SO_2)$-, phenyl-NH-$(SO_2)$-, $(\text{phenyl})_2$N-(SO_2)-, $(C_1\text{-}C_6)$alkyl-(C=O)-, phenyl-(C=O)-, $(C_3\text{-}C_{10})$cycloalkyl-(C=O)-, $(C_1\text{-}C_6)$alkyl-O-(C=O)-, $(C_3\text{-}C_{10})$cycloalkyl-O-(C=O)-, H_2N-(C=O)-, $(C_1\text{-}C_6)$alkyl-NH-(C=O)-, phenyl-NH-(C=O)-, $(C_3\text{-}C_{10})$cycloalkyl-NH-(C=O)-, $((C_1\text{-}C_6)\text{alkyl})_2$N-(C=O)-, $(\text{phenyl})_2$N-(C=O)-, phenyl-[($(C_1\text{-}C_6)$alkyl)-N]-(C=O)-,

and (C_3-C_{10})cycloalkyl-[((C_1-C_6)alkyl)-N]-(C=O)-;

where alkyl, alkenyl, alkynyl, phenyl, benzyl, cycloalkyl, alkoxy, phenoxy, amino of R^6 is optionally substituted with at least one moiety independently selected from the group consisting of halo, (C_1-C_6)alkyl,
(C_2-C_6)alkenyl, (C_2-C_6)alkynyl, perhalo(C_1-C_6)alkyl, (C_3-C_{10})cycloalkyl, phenyl, benzyl, ~~(C_5-C_{10})heterocyclic, (C_5-C_{10})heteroaryl~~, (C_1-C_6)alkyl-SO_2-, formyl, NC-,
(C_1-C_6)alkyl-(C=O)-, (C_3C_{10})cycloalkyl-(C=O)-, phenyl-(C=O)-,
~~(C_5-C_{10})heterocyclic-(C=O)-, (C_5-C_{10})heteroaryl-(C=O)-~~, HO-(C=O)-,
(C_1-C_6)alkyl-O-(C=O)-, (C_3-C_{10})cycloalkyl-O-(C=O)-,
~~(C_5-C_{10})heterocyclic-O-(C=O)-~~, (C_1-C_6)alkyl-NH-(C=O)-,
(C_3-C_{10})cycloalkyl-NH-(C=O)-, phenyl-NH-(C=O)-,
~~(C_5-C_{10})heterocyclic-NH-(C=O)-, (C_5-C_{10})heteroaryl-NH-(C=O)-~~,
((C_1-C_6)alkyl)$_2$-N-(C=O)-, phenyl-[((C_1-C_6)alkyl)-N]-(C=O)-, hydroxy,
(C_1-C_6)alkoxy, perhalo(C_1-C_6)alkoxy, (C_3-C_{10})cycloalkyl-O-, phenoxy,
~~(C_5-C_{10})heterocyclic-O-, (C_5-C_{10})heteroaryl-O-~~, (C_1-C_6)alkyl-(C=O)-O-,
(C_3-C_{10})cycloalkyl-(C=O)-O-, phenyl-(C=O)-O-, ~~(C_5-C_{10})heterocyclic-(C=O)-O-,~~
~~(C_5-C_{10})heteroaryl-(C=O)-O-~~, O_2N-, amino, (C_1-C_6)alkylamino,
((C_1-C_6)alkyl)$_2$-amino, formamidyl, (C_1-C_6)alkyl-(C=O)-NH-,
(C_3-C_{10})cycloalkyl-(C=O)-NH-, phenyl-(C=O)-NH-,
~~(C_5-C_{10})heterocyclic-(C=O)-NH-, (C_5-C_{10})heteroaryl-(C=O)-NH-~~,
(C_1-C_6)alkyl-(C=O)-[((C_1-C_6)alkyl)-N]-, phenyl-(C=O)-[(C_1-C_6)alkyl-N]-,
(C_1-C_6)alkyl-SO_2NH-, (C_3-C_{10})cycloalkyl-SO_2NH-, phenyl-SO_2NH-,
~~(C_5-C_{10})heterocyclic-SO_2NH- and (C_5-C_{10})heteroaryl-SO_2NH-~~;

wherein the phenyl moiety of a R^6 substituent is optionally further substituted with at least one radical independently selected from the group consisting of halo, (C_1-C_6)alkyl, (C_1-C_6)alkoxy, perfluoro(C_1-C_6)alkyl and perfluoro(C_1-C_6)alkoxy.

2. (Previously Cancelled)

3. (Previously Cancelled)

4. (Previously Cancelled)

5. (Previously Cancelled)

6. (Previously Cancelled)

7. (Previously Cancelled)

8. (Previously Cancelled)

9. (Original) A compound of claim 1, wherein s is one to two; R^3 is hydrogen or (C_1-C_6)alkyl; and R^6 is H, (C_1-C_6)alkyl, or (C_3-C_{10})cycloalkyl.

10. (Original) A pharmaceutical composition comprising a compound of claim 1 and a pharmaceutically acceptable carrier.

11. (Cancelled)

12. (Cancelled)

13. (Previously Presented) A compound 6-[5-(6-methyl-pyridin-2-yl)-2H-[1,2,3] triazol-4-yl]-quinazoline or a pharmaceutically acceptable salt thereof.

14. (Previously Presented) A pharmaceutical composition comprising 6-[5-(6-methyl-pyridin-2-yl)-2H-[1,2,3] triazol-4-yl]-quinazoline or a pharmaceutically acceptable salt thereof and a pharmaceutically acceptable carrier.